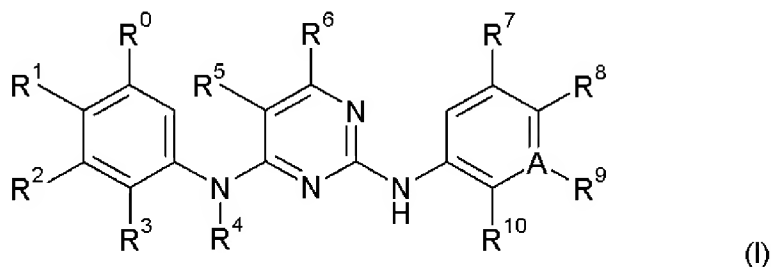


## Amendments to the Claims

This Listing of the Claims will replace all prior versions, and listings, of claims in the application.

### Listing of the Claims:

1. (Previously Presented) A compound of formula I



wherein

each of R<sup>0</sup>, R<sup>1</sup>, and R<sup>2</sup>, independently is hydrogen, C<sub>1</sub>-C<sub>8</sub>alkyl, C<sub>2</sub>-C<sub>8</sub>alkenyl, C<sub>2</sub>-C<sub>8</sub>alkinyl, C<sub>3</sub>-C<sub>8</sub>cycloalkyl, C<sub>3</sub>-C<sub>8</sub>cycloalkylC<sub>1</sub>-C<sub>8</sub>alkyl, C<sub>5</sub>-C<sub>10</sub>arylC<sub>1</sub>-C<sub>8</sub>alkyl, hydroxyC<sub>1</sub>-C<sub>8</sub>alkyl, C<sub>1</sub>-C<sub>8</sub>alkoxyC<sub>1</sub>-C<sub>8</sub>alkyl, aminoC<sub>1</sub>-C<sub>8</sub>alkyl, haloC<sub>1</sub>-C<sub>8</sub>alkyl, unsubstituted or substituted C<sub>5</sub>-C<sub>10</sub>aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1, 2 or 3 hetero atoms selected from N, O and S, hydroxy, C<sub>1</sub>-C<sub>8</sub>alkoxy, hydroxyC<sub>1</sub>-C<sub>8</sub>alkoxy, C<sub>1</sub>-C<sub>8</sub>alkoxyC<sub>1</sub>-C<sub>8</sub>alkoxy, haloC<sub>1</sub>-C<sub>8</sub>alkoxy, unsubstituted or substituted C<sub>5</sub>-C<sub>10</sub>arylC<sub>1</sub>-C<sub>8</sub>alkoxy, unsubstituted or substituted heterocyclyloxy, or unsubstituted or substituted heterocyclylC<sub>1</sub>-C<sub>8</sub>alkoxy, unsubstituted or substituted amino, C<sub>1</sub>-C<sub>8</sub>alkylthio, C<sub>1</sub>-C<sub>8</sub>alkylsulfinyl, C<sub>1</sub>-C<sub>8</sub>alkylsulfonyl, C<sub>5</sub>-C<sub>10</sub>arylsulfonyl, halogen, carboxy, C<sub>1</sub>-C<sub>8</sub>alkoxycarbonyl, unsubstituted or substituted carbamoyl, unsubstituted or substituted sulfamoyl, cyano or nitro;

R<sup>3</sup> is C<sub>1</sub>-C<sub>8</sub>alkylsulfinyl, C<sub>1</sub>-C<sub>8</sub>alkylsulfonyl, C<sub>5</sub>-C<sub>10</sub>arylsulfonyl, unsubstituted or substituted carbamoyl or unsubstituted or substituted sulfamoyl;

Or the pair of adjacent substituents R<sup>2</sup> and R<sup>3</sup> forms -CH<sub>2</sub>-NH-CO- or -CH<sub>2</sub>-NH-SO<sub>2</sub>- or such pairs wherein NH is substituted by C<sub>1</sub>-C<sub>8</sub>-alkyl;

R<sup>4</sup> is hydrogen or C<sub>1</sub>-C<sub>8</sub>alkyl;

each of R<sup>5</sup> and R<sup>6</sup> independently is hydrogen, C<sub>1</sub>-C<sub>8</sub>alkyl, C<sub>1</sub>-C<sub>8</sub>alkoxyC<sub>1</sub>-C<sub>8</sub>alkyl, haloC<sub>1</sub>-C<sub>8</sub>alkyl, C<sub>1</sub>-C<sub>8</sub>alkoxy, halogen, carboxy, C<sub>1</sub>-C<sub>8</sub>alkoxycarbonyl, unsubstituted or substituted carbamoyl, cyano, or nitro; and

each of R<sup>7</sup>, R<sup>8</sup>, R<sup>9</sup>, and R<sup>10</sup> independently is C<sub>1</sub>-C<sub>8</sub>alkyl, C<sub>2</sub>-C<sub>8</sub>alkenyl, C<sub>2</sub>-C<sub>8</sub>alkinyl, C<sub>3</sub>-C<sub>8</sub>cycloalkyl, C<sub>3</sub>-C<sub>8</sub>cycloalkylC<sub>1</sub>-C<sub>8</sub>alkyl, C<sub>5</sub>-C<sub>10</sub>arylC<sub>1</sub>-C<sub>8</sub>alkyl, hydroxyC<sub>1</sub>-C<sub>8</sub>alkyl, C<sub>1</sub>-

C<sub>8</sub>alkoxyC<sub>1</sub>-C<sub>8</sub>alkyl, aminoC<sub>1</sub>-C<sub>8</sub>alkyl, haloC<sub>1</sub>-C<sub>8</sub>alkyl, unsubstituted or substituted C<sub>5</sub>-C<sub>10</sub>aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1, 2 or 3 hetero atoms selected from N, O and S, hydroxy, C<sub>1</sub>-C<sub>8</sub>alkoxy, hydroxyC<sub>1</sub>-C<sub>8</sub>alkoxy, C<sub>1</sub>-C<sub>8</sub>alkoxyC<sub>1</sub>-C<sub>8</sub>alkoxy, haloC<sub>1</sub>-C<sub>8</sub>alkoxy, unsubstituted or substituted C<sub>5</sub>-C<sub>10</sub>arylC<sub>1</sub>-C<sub>8</sub>alkoxy, unsubstituted or substituted heterocyclyoxy, or unsubstituted or substituted heterocyclylC<sub>1</sub>-C<sub>8</sub>alkoxy, unsubstituted or substituted amino, C<sub>1</sub>-C<sub>8</sub>alkylthio, C<sub>1</sub>-C<sub>8</sub>alkylsulfinyl, C<sub>1</sub>-C<sub>8</sub>alkylsulfonyl, C<sub>5</sub>-C<sub>10</sub>arylsulfonyl, halogen, carboxy, C<sub>1</sub>-C<sub>8</sub>alkoxycarbonyl, unsubstituted or substituted carbamoyl, unsubstituted or substituted sulfamoyl, cyano or nitro; wherein R<sup>7</sup>, R<sup>8</sup> and R<sup>9</sup> independently of each other can also be hydrogen;

or R<sup>7</sup> and R<sup>8</sup>, R<sup>8</sup> and R<sup>9</sup>, and/or R<sup>9</sup> and R<sup>10</sup> form together with the carbon atoms to which they are attached, a 5 or 6 membered carbocyclic or heterocyclic ring comprising 0, 1, 2 or 3 heteroatoms selected from N, O and S that is unsubstituted or substituted by C<sub>1</sub>-C<sub>8</sub>alkyl, C<sub>1</sub>-C<sub>8</sub>alkoxy, halo-C<sub>1</sub>-C<sub>8</sub>alkyl, hydroxyl, amino, substituted amino, halogen, carboxy, C<sub>1</sub>-C<sub>8</sub>alkoxycarbonyl, carbamoyl, cyano, or oxo;

A is C;

and salts thereof.

2. (Currently Amended) A compound of formula I according to claim 1, wherein each of R<sup>0</sup> or R<sup>2</sup> independently is hydrogen, C<sub>1</sub>-C<sub>8</sub>alkyl, hydroxyC<sub>1</sub>-C<sub>8</sub>alkyl, haloC<sub>1</sub>-C<sub>8</sub>alkyl, unsubstituted or substituted C<sub>5</sub>-C<sub>10</sub>aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S, C<sub>1</sub>-C<sub>8</sub>alkoxy, haloC<sub>1</sub>-C<sub>8</sub>alkoxy, C<sub>5</sub>-C<sub>10</sub>aryloxy, unsubstituted or substituted heterocyclyoxy, unsubstituted or substituted heterocyclylC<sub>1</sub>-C<sub>8</sub>alkoxy, unsubstituted or substituted amino, C<sub>1</sub>-C<sub>8</sub>alkylsulfonyl, halogen, unsubstituted or substituted carbamoyl, unsubstituted or substituted sulfamoyl;

R<sup>1</sup> is hydrogen, C<sub>1</sub>-C<sub>8</sub>alkyl, hydroxyC<sub>1</sub>-C<sub>8</sub>alkyl, haloC<sub>1</sub>-C<sub>8</sub>alkyl, unsubstituted or substituted C<sub>5</sub>-C<sub>10</sub>aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S, C<sub>1</sub>-C<sub>8</sub>alkoxy, haloC<sub>1</sub>-C<sub>8</sub>alkoxy, C<sub>5</sub>-C<sub>10</sub>aryloxy, unsubstituted or substituted heterocyclyoxy, unsubstituted or substituted heterocyclylC<sub>1</sub>-C<sub>8</sub>alkoxy, unsubstituted or substituted amino, C<sub>1</sub>-C<sub>8</sub>alkylsulfonyl, halogen, unsubstituted or substituted carbamoyl, unsubstituted or substituted sulfamoyl;

R<sup>3</sup> is C<sub>1</sub>-C<sub>8</sub>alkylsulfinyl, C<sub>1</sub>-C<sub>8</sub>alkylsulfonyl, C<sub>5</sub>-C<sub>10</sub>arylsulfonyl, unsubstituted or substituted carbamoyl or unsubstituted or substituted sulfamoyl;

~~R<sup>3</sup> is C<sub>1</sub>-C<sub>8</sub>alkylsulfinyl, C<sub>1</sub>-C<sub>8</sub>alkylsulfonyl, C<sub>5</sub>-C<sub>10</sub>arylsulfonyl, unsubstituted or substituted carbamoyl or unsubstituted or substituted sulfamoyl;~~

or the pair of adjacent substituents  $R^2$  and  $R^3$  forms  $-CH_2-NH-CO-$  or  $CH_2-NH-SO_2-$  or such pairs wherein NH is substituted by  $C_1-C_8$ -alkyl;

$R^4$  is hydrogen or  $C_1-C_8$ -alkyl;

$R^5$  is hydrogen;  $C_1-C_8$ -alkyl, halogen, halo $C_1-C_8$ -alkyl, cyano or nitro;

$R^6$  is hydrogen;

each of  $R^7$  and  $R^9$  independently is hydrogen,  $C_1-C_8$ -alkyl, hydroxy $C_1-C_8$ -alkyl, halo $C_1-C_8$ -alkyl, unsubstituted or substituted  $C_5-C_{10}$ -aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S,  $C_1-C_8$ -alkoxy, halo $C_1-C_8$ -alkoxy,  $C_5-C_{10}$ -aryloxy, unsubstituted or substituted heterocyclyloxy, unsubstituted or substituted heterocyclyl $C_1-C_8$ -alkoxy, unsubstituted or substituted amino,  $C_1-C_8$ -alkylsulfonyl, halogen, unsubstituted or substituted carbamoyl, unsubstituted or substituted sulfamoyl;

$R^8$  is hydrogen,  $C_1-C_8$ -alkyl, hydroxy $C_1-C_8$ -alkyl, halo $C_1-C_8$ -alkyl,  $C_5-C_{10}$ -aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S,  $C_1-C_8$ -alkoxy, halo $C_1-C_8$ -alkoxy,  $C_5-C_{10}$ -aryloxy, unsubstituted or substituted heterocyclyloxy, unsubstituted or substituted heterocyclyl $C_1-C_8$ -alkoxy, unsubstituted or substituted amino,  $C_1-C_8$ -alkylsulfonyl, halogen, unsubstituted or substituted carbamoyl, unsubstituted or substituted sulfamoyl, cyano, or nitro; and

$R^{10}$  is  $C_1-C_8$ -alkyl, hydroxy $C_1-C_8$ -alkyl, halo $C_1-C_8$ -alkyl,  $C_1-C_8$ -alkoxy, unsubstituted or substituted heterocyclyl $C_1-C_8$ -alkoxy, unsubstituted or substituted amino, halogen, carboxy, carbamoyl, or unsubstituted or substituted sulfamoyl; or

each pair of adjacent substituents  $R^7$  and  $R^8$ , or  $R^8$  and  $R^9$  or  $R^9$  and  $R^{10}$ , is  $-NH-CH=CH-$ ,  $-CH=CH-NH-$ ,  $-NH-N=CH-$ ,  $-CH=N-NH-$ ,  $-CH_2-CH_2-CH_2-$ ,  $-CH_2-CH_2-CH_2-CH_2-$ ,  $-CH_2-CH_2-O-$ ,  $-CH=CH-O-$ ,  $-O-CH_2-O-$ , or  $-O-CF_2-O-$ ;

A is C.

3. (Previously Presented) A compound of formula I according to claim 1, wherein

each of  $R^0$  or  $R^2$  independently is hydrogen,  $C_1-C_8$ -alkyl, halo $C_1-C_8$ -alkyl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S,  $C_1-C_8$ -alkoxy, unsubstituted or substituted heterocyclyloxy, unsubstituted or substituted heterocyclyl $C_1-C_8$ -alkoxy, unsubstituted or substituted amino, or halogen;

$R^1$  is hydrogen,  $C_1-C_8$ -alkyl, halo $C_1-C_8$ -alkyl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S,  $C_1-C_8$ -alkoxy, unsubstituted or substituted heterocyclyloxy, unsubstituted or substituted heterocyclyl $C_1-C_8$ -alkoxy, unsubstituted or substituted amino, halogen;

$R^3$  is  $C_1$ - $C_8$ alkylsulfinyl,  $C_1$ - $C_8$ -alkylsulfonyl,  $C_5$ - $C_{10}$ arylsulfonyl, unsubstituted or substituted carbamoyl or unsubstituted or substituted sulfamoyl;  
 or the pair of adjacent substituents  $R^2$  and  $R^3$  forms  $-CH_2-NH-CO-$  or  $-CH_2-NH-SO_2-$  or such pairs wherein NH is substituted by  $C_1$ - $C_8$ -alkyl;  
 $R^4$  is hydrogen;  
 $R^5$  is hydrogen, halogen, halo $C_1$ - $C_8$ alkyl, or nitro;  
 $R^6$  is hydrogen;  
 each of  $R^7$  and  $R^9$  independently is hydrogen,  $C_1$ - $C_8$ alkyl, halo $C_1$ - $C_8$ alkyl, unsubstituted or substituted  $C_5$ - $C_{10}$ aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S,  $C_1$ - $C_8$ alkoxy, unsubstituted or substituted heterocycloxy, unsubstituted or substituted heterocyclyl $C_1$ - $C_8$ alkoxy, unsubstituted or substituted amino, halogen, unsubstituted or substituted carbamoyl, or unsubstituted or substituted sulfamoyl;  
 $R^8$  is hydrogen,  $C_1$ - $C_8$ alkyl, halo $C_1$ - $C_8$ alkyl,  $C_5$ - $C_{10}$ aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S,  $C_1$ - $C_8$ alkoxy, halo $C_1$ - $C_8$ alkoxy,  $C_5$ - $C_{10}$ aryloxy, unsubstituted or substituted heterocycloxy, unsubstituted or substituted heterocyclyl $C_1$ - $C_8$ alkoxy, unsubstituted or substituted amino, halogen, unsubstituted or substituted sulfamoyl, or nitro; and  
 $R^{10}$  is  $C_1$ - $C_8$ alkyl, halo $C_1$ - $C_8$ alkyl,  $C_1$ - $C_8$ alkoxy, unsubstituted or substituted heterocyclyl $C_1$ - $C_8$ alkoxy, unsubstituted or substituted amino, or halogen; or  
 each pair of adjacent substituents  $R^7$  and  $R^8$ , or  $R^8$  and  $R^9$  or  $R^9$  and  $R^{10}$ , is  $-NH-CH=CH-$ ,  $-CH=CH-NH-$ ,  $-NH-N=CH-$ ,  $-CH=N-NH-$ ,  $-CH_2-CH_2-CH_2-$ ,  $-CH_2-CH_2-CH_2-CH_2-$ ,  $-O-CH_2-O-$ , or  $-O-CF_2-O-$ ;  
 A is C.

4. (Currently Amended) A compound of formula I according to claim 1, wherein  
 each of  $R^0$  or  $R^2$  independently is hydrogen, piperazino, N-methylpiperazino or 1-methyl-4-piperidyloxy;  
 $R^1$  is hydrogen, piperazino, N-methylpiperazino, morpholino, 1-methyl-4-piperidinyloxy, 3-morpholinopropoxy or 2-morpholinoethoxy;  
 $R^3$  is sulfamoyl, methylsulfamoyl or propylsulfamoyl; ~~or~~  
~~the pair of adjacent substituents  $R^0$  and  $R^1$ , or  $R^1$  and  $R^2$  is  $-O-CH_2-O-$ , or the pair of~~  
 adjacent substituents  $R^2$  and  $R^3$  is  $-CH_2-NH-CO-$  or  $-CH_2-NH-SO_2-$ ;  
 $R^4$  is hydrogen;  
 $R^5$  is hydrogen, chloro, bromo, trifluoromethyl or nitro;  
 $R^6$  is hydrogen;

each of  $R^7$  and  $R^9$  independently is hydrogen,  $C_1$ - $C_8$ alkyl, halo $C_1$ - $C_8$ alkyl, unsubstituted or substituted  $C_5$ - $C_{10}$ aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S,  $C_1$ - $C_8$ alkoxy, unsubstituted or substituted heterocyclyloxy, unsubstituted or substituted heterocyclyl $C_1$ - $C_8$ alkoxy, unsubstituted or substituted amino, halogen, unsubstituted or substituted carbamoyl, or unsubstituted or substituted sulfamoyl;

$R^8$  is hydrogen,  $C_1$ - $C_8$ alkyl, halo $C_1$ - $C_8$ alkyl,  $C_5$ - $C_{10}$ aryl, unsubstituted or substituted 5 or 6 membered heterocyclyl comprising 1 or 2 hetero atoms selected from N, O and S,  $C_1$ - $C_8$ alkoxy, halo $C_1$ - $C_8$ alkoxy,  $C_5$ - $C_{10}$ aryloxy, unsubstituted or substituted heterocyclyloxy, unsubstituted or substituted heterocyclyl $C_1$ - $C_8$ alkoxy, unsubstituted or substituted amino, halogen, unsubstituted or substituted sulfamoyl, or nitro; and

$R^{10}$  is  $C_1$ - $C_8$ alkyl, halo $C_1$ - $C_8$ alkyl,  $C_1$ - $C_8$ alkoxy, unsubstituted or substituted heterocyclyl $C_1$ - $C_8$ alkoxy, unsubstituted or substituted amino, or halogen; or

each pair of adjacent substituents  $R^7$  and  $R^8$ , or  $R^8$  and  $R^9$  or  $R^9$  and  $R^{10}$ , is  $-NH-CH=CH-$ ,  $-CH=CH-NH-$ ,  $-NH-N=CH-$ ,  $-CH=N-NH-$ ,  $-CH_2-CH_2-CH_2-$ ,  $-CH_2-CH_2-CH_2-CH_2-$ ,  $-O-CH_2-O-$ , or  $-O-CF_2-O-$ ;

A is C.

5. (Currently Amended) A compound of formula I according to claim 1, wherein each of  $R^0$  or  $R^2$  independently is hydrogen, piperazino, N-methylpiperazino or 1-methyl-4-piperidyloxy;

$R^1$  is hydrogen, piperazino, N-methylpiperazino, morpholino, 1-methyl-4-piperidinyloxy, 3-morpholinopropoxy or 2-morpholinoethoxy;

$R^3$  is sulfamoyl, methylsulfamoyl or propylsulfamoyl; or

~~the pair of adjacent substituents  $R^0$  and  $R^1$ , or  $R^1$  and  $R^2$  is  $-O-CH_2-O-$ , or the pair of adjacent substituents  $R^2$  and  $R^3$  is  $-CH_2-NH-CO-$  or  $-CH_2-NH-SO_2-$ ;~~

$R^4$  is hydrogen;

$R^5$  is hydrogen, chloro, bromo, trifluoromethyl or nitro;

$R^6$  is hydrogen;

each of  $R^7$  and  $R^9$  independently is hydrogen, methyl, isopropyl, trifluoromethyl, phenyl, o-, m- or p-methoxyphenyl, piperidino, piperazino, N-methylpiperazino, morpholino, methoxy, ethoxy, isopropoxy, phenoxy, 3-morpholinopropoxy, 2-morpholinoethoxy, 2-(1-imidazolyl)ethoxy, dimethylamino, fluoro, morpholinocarbonyl, piperidinocarbonyl, piperazinocarbonyl or cyclohexylcarbamoyl;

$R^8$  is hydrogen, methyl, piperidino, piperazino, N-methylpiperazino, morpholino, methoxy, ethoxy, trifluoromethoxy, phenoxy, 1-methyl-4-piperidyloxy, 3-morpholinopropoxy, 2-

morpholinoethoxy, 3-(N-methylpiperazino)-propoxy, methylamino, fluoro, chloro, sulfamoyl or nitro; and

R<sup>10</sup> is methyl, butyl, methoxy, ethoxy, 2-(1-imidazolyl)ethoxy, methylamino, dimethylamino or fluoro; or

the pair of adjacent substituents R<sup>7</sup> and R<sup>8</sup> or R<sup>8</sup> and R<sup>9</sup> is -O-CH<sub>2</sub>-O- or the pair of adjacent substituents R<sup>9</sup> and R<sup>10</sup> is -NH-CH=CH-, -CH=N-NH-, -CH<sub>2</sub>-CH<sub>2</sub>-CH<sub>2</sub>-, -CH<sub>2</sub>-CH<sub>2</sub>-CH<sub>2</sub>-CH<sub>2</sub>- or -O-CF<sub>2</sub>-O-;

A is C.

6. (Previously Presented) A compound of formula I according to claim 1, wherein each of R<sup>0</sup>, R<sup>1</sup> or R<sup>2</sup> is hydrogen;

R<sup>3</sup> is sulfamoyl, methylsulfamoyl or propylsulfamoyl;

R<sup>4</sup> is hydrogen;

R<sup>5</sup> is chloro or bromo;

R<sup>6</sup> is hydrogen;

each of R<sup>7</sup> and R<sup>9</sup> independently is hydrogen, methyl, isopropyl, trifluoromethyl, phenyl, o-, m- or p-methoxyphenyl, piperidino, piperazino, N-methylpiperazino, morpholino, methoxy, ethoxy, isopropoxy, phenoxy, 3-morpholinopropoxy, 2-morpholinoethoxy, 2-(1-imidazolyl)ethoxy, dimethylamino, fluoro, morpholinocarbonyl, piperidinocarbonyl, piperazinocarbonyl or cyclohexylcarbonyl;

R<sup>8</sup> is hydrogen, methyl, piperidino, piperazino, N-methylpiperazino, morpholino, methoxy, ethoxy, trifluoromethoxy, phenoxy, 1-methyl-4-piperidyloxy, 3-morpholinopropoxy, 2-morpholinoethoxy, 3-(N-methylpiperazino)-propoxy, methylamino, fluoro, chloro, sulfamoyl or nitro; and

R<sup>10</sup> is methyl, butyl, methoxy, ethoxy, 2-(1-imidazolyl)ethoxy, methylamino, dimethylamino or fluoro; or

the pair of adjacent substituents R<sup>7</sup> and R<sup>8</sup> or R<sup>8</sup> and R<sup>9</sup> is -O-CH<sub>2</sub>-O-, or the pair of adjacent substituents R<sup>9</sup> and R<sup>10</sup> is -NH-CH=CH-, -CH=N-NH-, -CH<sub>2</sub>-CH<sub>2</sub>-CH<sub>2</sub>-, -CH<sub>2</sub>-CH<sub>2</sub>-CH<sub>2</sub>-CH<sub>2</sub>- or -O-CF<sub>2</sub>-O-;

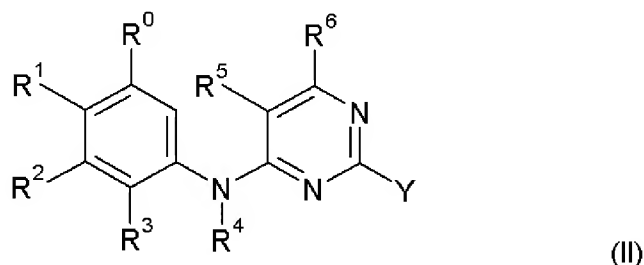
A is C.

7. (Previously Presented) The compound of formula I according to claim 1, wherein each of R<sup>0</sup>, R<sup>1</sup> or R<sup>2</sup> is hydrogen, R<sup>3</sup> is methylsulfamoyl, R<sup>4</sup> is hydrogen, R<sup>5</sup> is bromo, R<sup>6</sup> is hydrogen, each of R<sup>7</sup> and R<sup>8</sup> is methoxy, R<sup>9</sup> is hydrogen, and R<sup>10</sup> is methyl, and A is C.

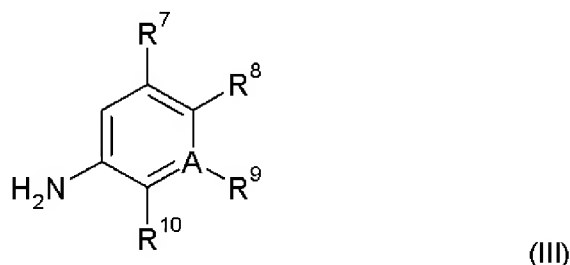
8. (Previously Presented) The compound of formula I according to claim 1, wherein each of  $R^0$ ,  $R^1$  or  $R^2$  is hydrogen,  $R^3$  is methylsulfamoyl,  $R^4$  is hydrogen,  $R^5$  is bromo,  $R^6$  is hydrogen, each of  $R^7$  and  $R^8$  is hydrogen, and the pair of adjacent substituents  $R^9$  and  $R^{10}$  is  $-\text{CH}_2-\text{CH}_2-\text{CH}_2-$ , and A is C.

9. (Previously Presented) The compound 2-{5-Chloro-2-[4-(3-methylamino-pyrrolidin-1-yl)-phenylamino]-pyrimidin-4-ylamino}-N-isopropyl-benzenesulfonamide.

10. (Previously Presented) A process for the production of a compound of formula I according to claim 1, comprising reacting a compound of formula II



wherein  $R^0$ ,  $R^1$ ,  $R^2$ ,  $R^3$ ,  $R^4$ ,  $R^5$ , and  $R^6$  are as defined in claim 1, and Y is a leaving group, with a compound of formula III



wherein A,  $R^7$ ,  $R^8$ ,  $R^9$  and  $R^{10}$  are as defined in claim 1;

and, if desired, converting a compound of formula I, wherein the substituents have the meaning as defined in claim 1, into another compound of formula I as defined in claim 1;

and recovering the resulting compound of formula I in free form or as a salt, and, when required, converting the compound of formula I obtained in free form into the desired salt, or an obtained salt into the free form.

11. (Previously Presented) A pharmaceutical composition comprising a compound according to claim 1, as active ingredient together with one or more pharmaceutically acceptable diluents or carriers.

12. (Cancelled).

13. (Previously Presented) A combination comprising a therapeutically effective amount of a compound according to claim 1 and one or more known drug substances, said further drug substance being useful in the treatment of neoplastic diseases or immune system disorders.

14. (Currently Amended) A method for the treatment of breast tumors ~~neoplastic diseases and immune system disorders~~ in a subject in need thereof which comprises administering an effective amount of a compound according to claim 1 or a pharmaceutical composition comprising same.

15. (Currently Amended) A method for the treatment ~~or prevention~~ of a disease which responds to inhibition of focal adhesion kinase or/and IGF-1 Receptor comprising administering a compound according to claim 1 or a pharmaceutically acceptable salt thereof.

16.-20. (Cancelled).

21. (Previously Presented) The method according to claim 14, wherein the compound is 2-[5-Bromo-2-(2-methoxy-5-morpholin-4-yl-phenylamino)-pyrimidin-4-ylamino]-N-methyl-benzenesulfonamide or a pharmaceutically acceptable salt thereof.

22. (Previously Presented) The method according to claim 14, wherein the compound is selected from 2-[5-chloro-2-(2-methoxy-4-morpholin-4-yl-phenylamino)-pyrimidin-4-ylamino]-N-methyl-benzamide, N<sup>2</sup>-(4-[1,4']Bipiperidiny-1'-yl-2-methoxy-phenyl)-5-chloro-N<sup>4</sup>-[2-(propane-1-sulfonyl)-phenyl]-pyrimidine-2,4-diamine and 2-{5-Chloro-2-[2-methoxy-4-(4-methyl-piperazin-1-yl)-phenylamino]-pyrimidin-4-ylamino}-N-isopropyl-benzenesulfonamide, or a pharmaceutically acceptable salt thereof.

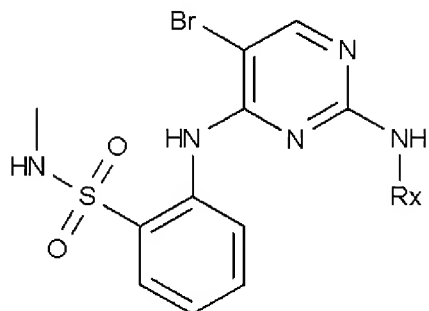
23. (Previously Presented) A compound of the formula I shown in claim 1, selected from the group of compounds with the following names or formulae:

2-[2-(2,5-dimethoxy-phenylamino)-5-nitro-pyrimidin-4-ylamino]-N-methyl-benzenesulfonamide;

2-[5-bromo-2-(2,4-dimethoxy-phenylamino)-pyrimidin-4-ylamino]-N-methyl-benzenesulfonamide;

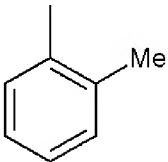
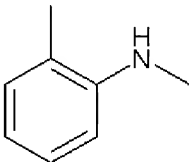
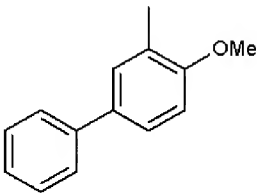
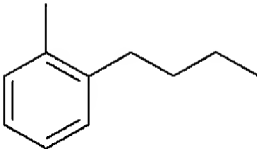
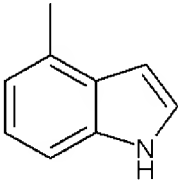
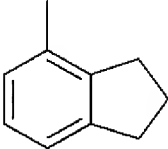
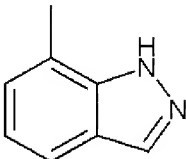
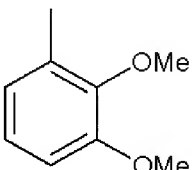


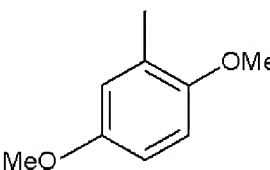
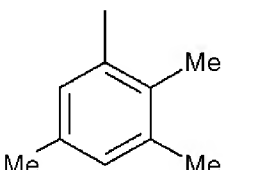
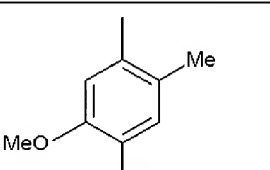
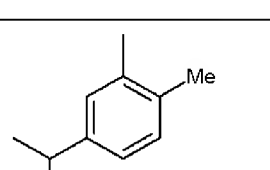
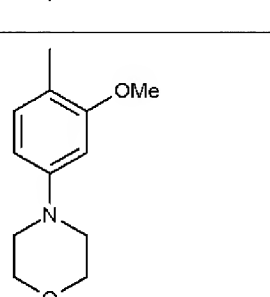
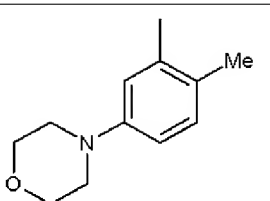
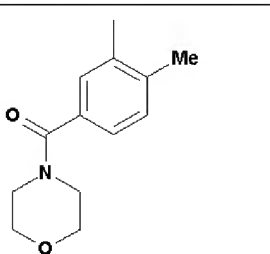
a compound of the formula

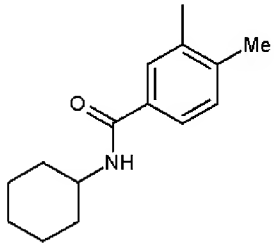
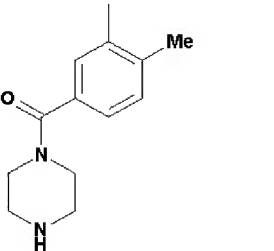
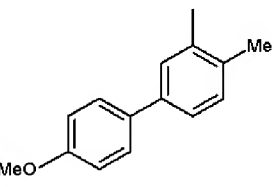
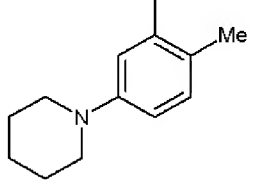
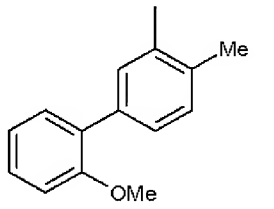
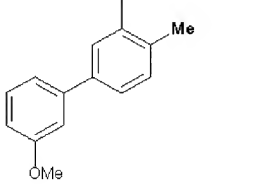
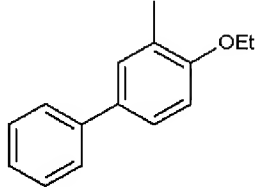


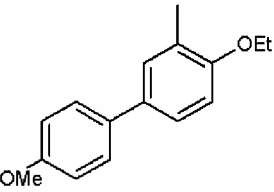
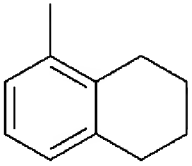
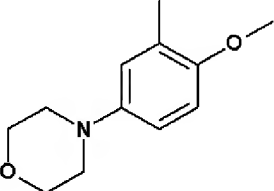
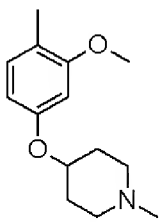
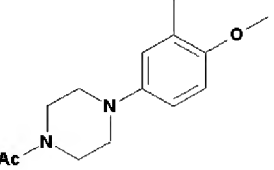
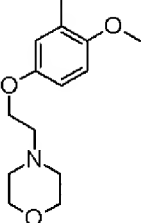
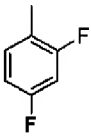
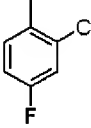
wherein Rx has one of the meanings given in the following table:

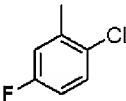
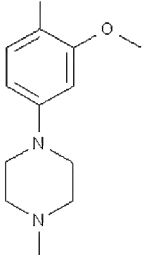
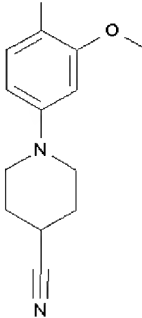
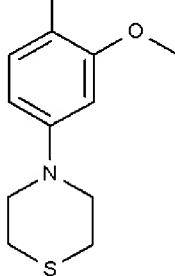
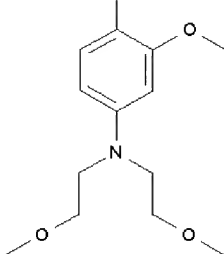
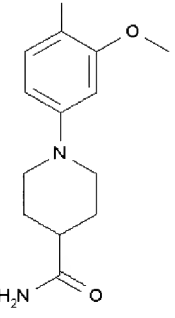
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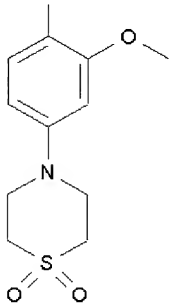
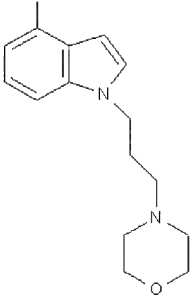
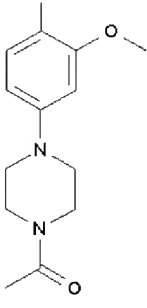
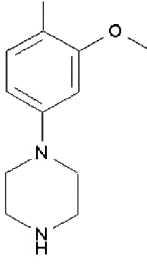
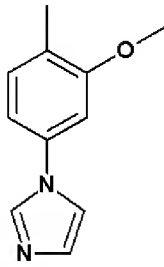
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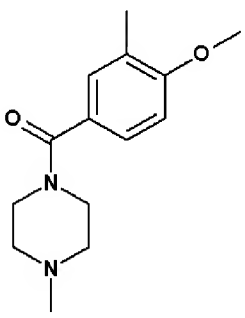
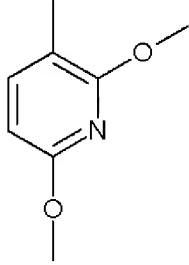
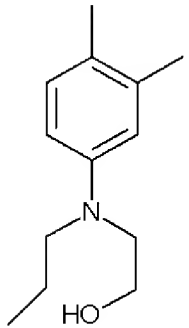
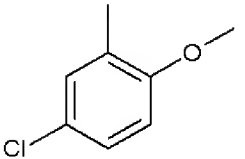
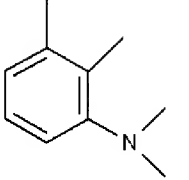
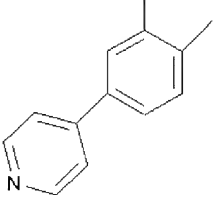
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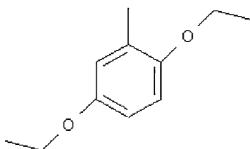
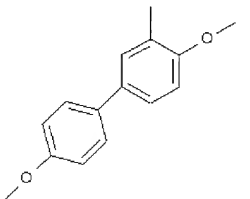
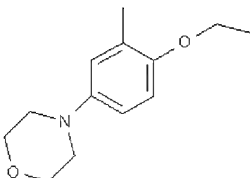
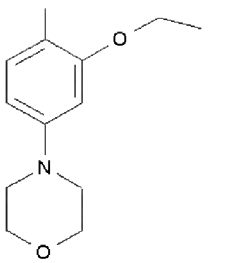
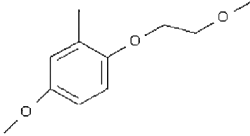
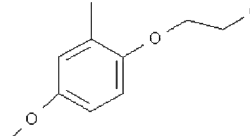
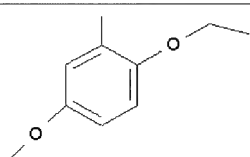
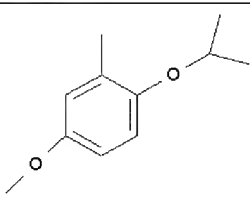
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3-31	 <chem>CC1=CC=C2CCCCC12</chem>
3-32	 <chem>COc1ccc(cc1)N2CCCC2</chem>
3-33	 <chem>COc1ccc(cc1)N2CCN(C)CC2</chem>
3-34	 <chem>CC(=O)N1CCN(Cc2ccc(OC)cc2)CC1</chem>
3-35	 <chem>COc1ccc(cc1)OCCN2CCOCC2</chem>
3-36	 <chem>Fc1cccc(F)c1</chem>
3-37	 <chem>Fc1cccc(Cl)c1</chem>

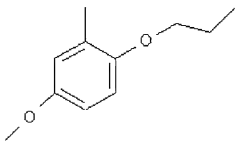
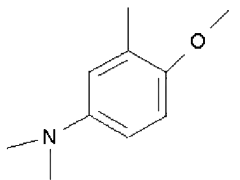
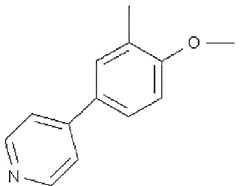
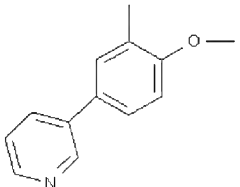
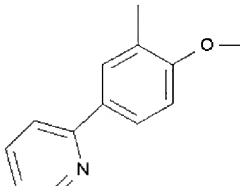
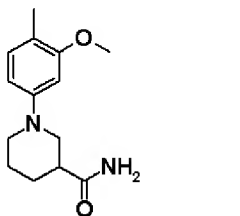
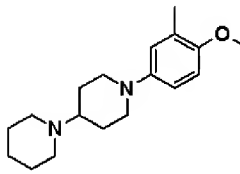
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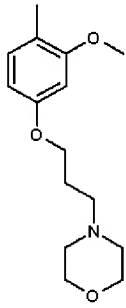
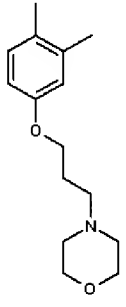
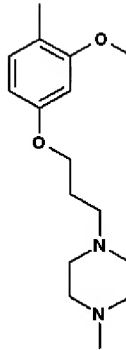
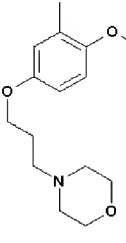
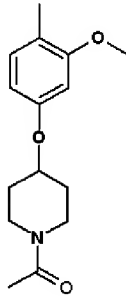
3-44	 <chem>COC1=CC=C(C=C1)N2CCSC(=O)(=O)CC2</chem>
3-45	 <chem>CC1=CC=C2C(=C1)C(=CN2)CCCN3CCOCC3</chem>
3-46	 <chem>CC(=O)N1CCN(CC1)C2=CC=C(OC)C=C2</chem>
3-47	 <chem>COC1=CC=C(C=C1)N2CCNCC2</chem>
3-48	 <chem>COC1=CC=C(C=C1)N2C=CC=C2</chem>

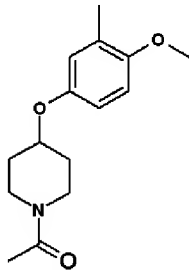
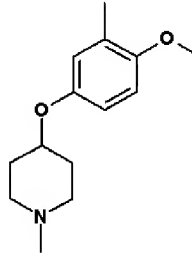
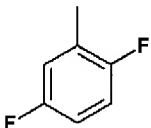
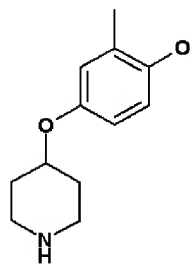
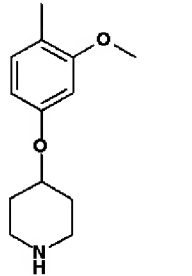
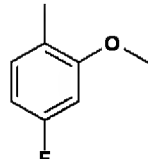
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3-50	 <chem>COC1=CC=CC(=O)N=C1OC</chem>
3-51	 <chem>CC1=CC=C(C=C1)N2CCCN2C3=CC=CC=C3C</chem>
3-52	 <chem>COC1=CC=C(C=C1)C(=O)N2CCCN2C3=CC=CC=C3C</chem>
3-53	 <chem>CC1=CC=C(C=C1)C(=O)N2CCCN2C3=CC=CC=C3C</chem>
3-54	 <chem>C1=CC=C(C=C1)C(=O)N2CCCN2C3=CC=CC=C3C</chem>



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3-70	 <chem>COc1cccc(c1)OCCCN2CCOCC2</chem>
3-71	 <chem>Cc1cc(C)ccc1OCCCN2CCOCC2</chem>
3-72	 <chem>COc1cccc(c1)OCCCN2CCN(C)CCN2C</chem>
3-73	 <chem>COc1cccc(c1)OCCCN2CCOCC2</chem>
3-74	 <chem>CC(=O)N1CCOC(c2ccc(OC)cc2)C1</chem>

3-75	 <chem>CC(=O)N1CCCC1Oc2cc(C)c(OC)cc2</chem>
3-76	 <chem>CN1CCCC1Oc2cc(C)c(OC)cc2</chem>
3-77	 <chem>Fc1cccc(F)c1</chem>
3-78	 <chem>CN1CCCC1Oc2cc(C)c(OC)cc2</chem>
3-79	 <chem>CN1CCCCC1Oc2cc(C)c(OC)cc2</chem>
3-80	 <chem>COc1cccc(F)c1</chem>

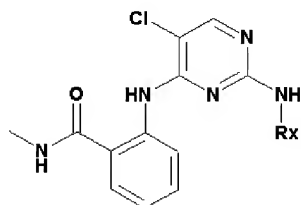
3-81	
3-82	
3-83	
3-84	

;

2-[5-bromo-2-(2,3-[difluoromethylenedioxy]phenylamino)-pyrimidin-4-ylamino]-benzenesulfonamide;

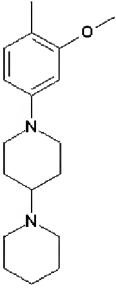
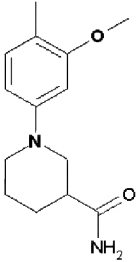
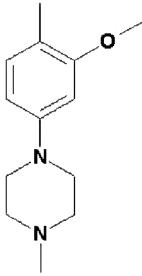
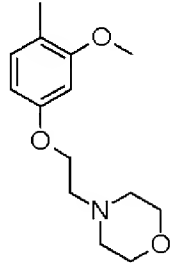
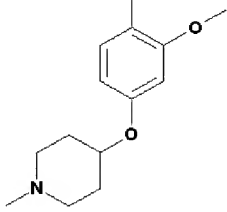
2-[5-chloro-2-(2-methoxy-4-morpholin-4-yl-phenylamino)-pyrimidin-4-ylamino]-N-methylbenzamide;

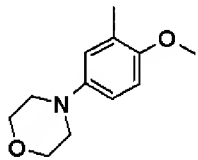
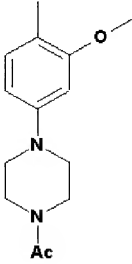
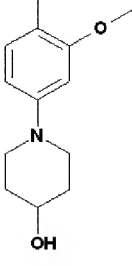
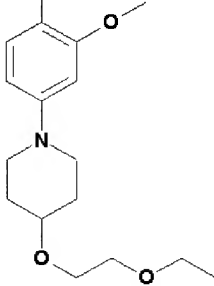
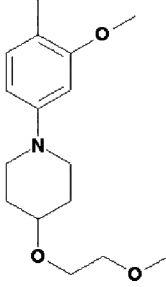
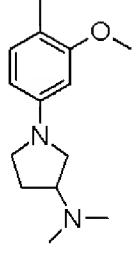
a compound of the formula

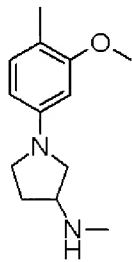
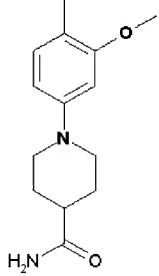
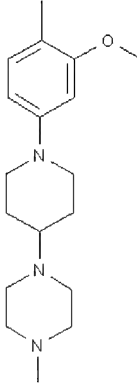
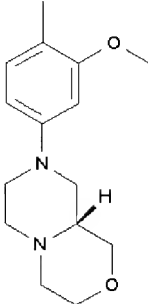
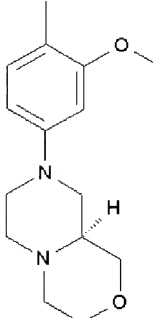


wherein Rx has one of the meanings given in the following table:

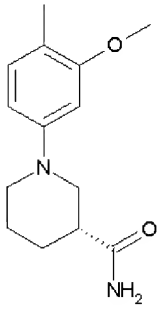
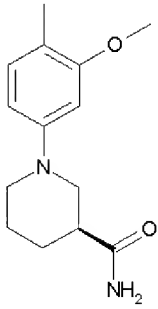
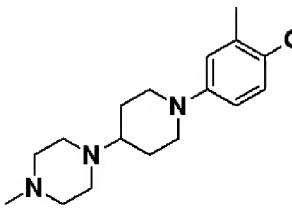
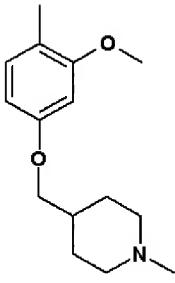
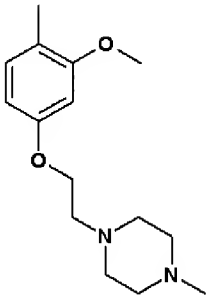
Compound	Rx

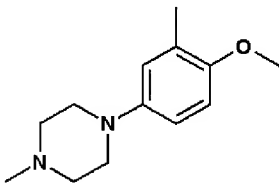
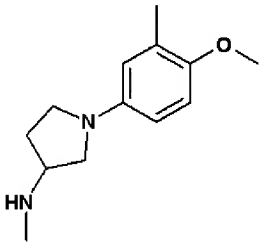
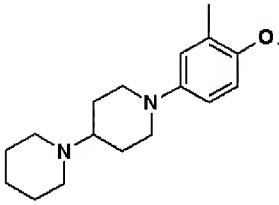
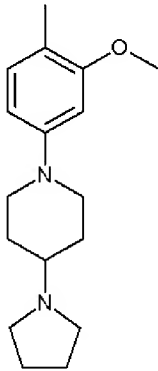
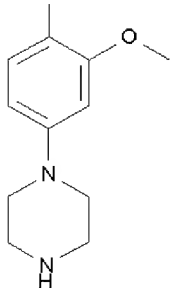
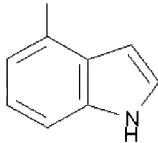
No.	
7-1	 <chem>COC1=CC=C(C=C1)N2CCN(CC2)C3CCCCN3</chem>
7-2	 <chem>COC1=CC=C(C=C1)N2CCN(CC2)C(=O)N</chem>
7-3	 <chem>COC1=CC=C(C=C1)N2CCN(CC2)</chem>
7-4	 <chem>COC1=CC=C(C=C1)OCCN2CCOCC2</chem>
7-5	 <chem>COC1=CC=C(C=C1)OCC2CCN(C)CC2</chem>

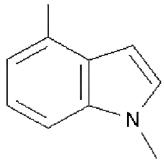
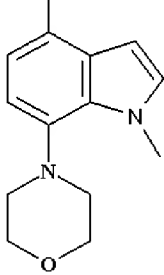
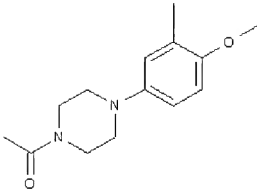
7-6	
7-7	
7-8	
7-9	
7-10	
7-11	

7-12	 <chem>CNCC1CN(C1)c2ccc(OC)cc2</chem>
7-13	 <chem>NC(=O)C1CCN(C1)c2ccc(OC)cc2</chem>
7-14	 <chem>CN1CCN(C1)C2CCN(C2)c3ccc(OC)cc3</chem>
7-15	 <chem>C1CCN(C1)C2CCN(C2)c3ccc(OC)cc3</chem>
7-16	 <chem>C1CCN(C1)C2CCN(C2)c3ccc(OC)cc3</chem>



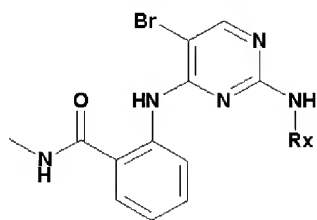
7-17	
7-18	
7-19	
7-20	
7-21	

7-22	
7-23	
7-24	
7-25	
7-26	
7-27	

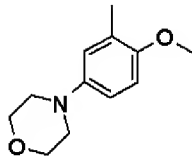
7-28	
7-29	
7-30	

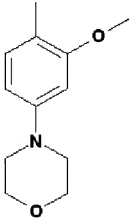
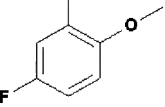
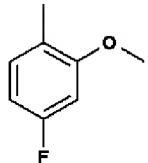
;

a compound of the formula



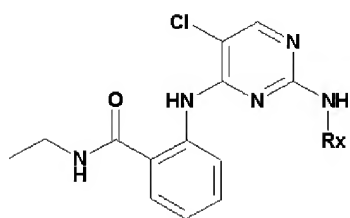
wherein Rx has one of the meanings given in the following table:

Com- pound	Rx
8-1	

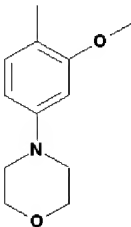
8-2	
8-3	
8-4	

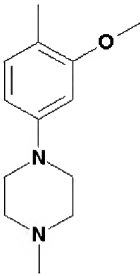
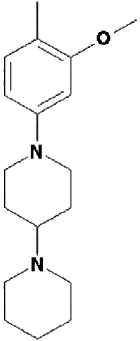
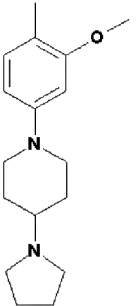
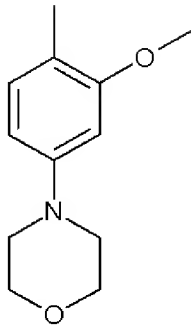
;

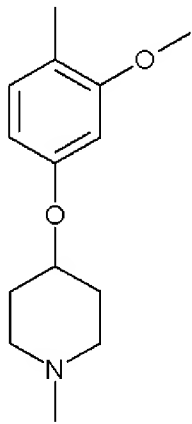
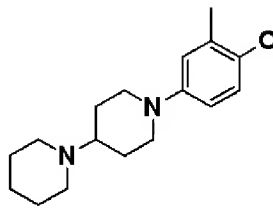
a compound of the formula



wherein Rx has one of the meanings given in the following table:

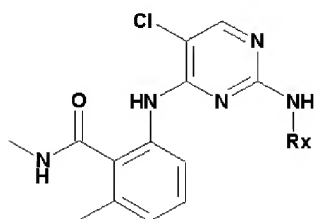
Com- pound	Rx
9-1	

9-2	 <chem>COC1=CC=C(C=C1)N2CCNCC2</chem>
9-3	 <chem>COC1=CC=C(C=C1)N2CCCCC2</chem>
9-4	 <chem>COC1=CC=C(C=C1)N2CCCC2</chem>
9-5	 <chem>COC1=CC=C(C=C1)N2CCOCC2</chem>

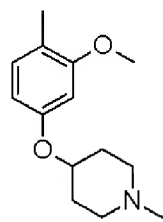
9-6	
9-7	

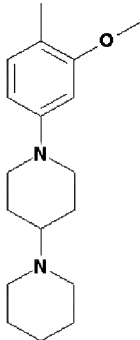
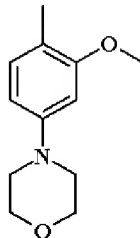
i

a compound of the formula



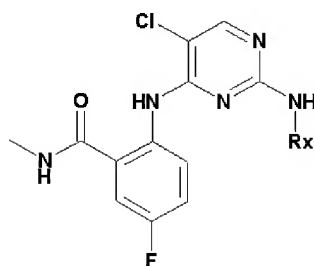
wherein Rx has one of the meanings given in the following table:

Com- pound	Rx
10-1	

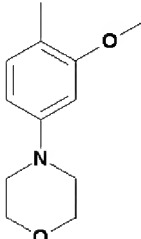
10-2	
10-3	

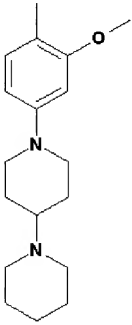
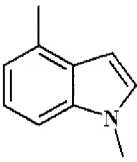
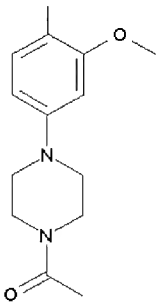
;

a compound of the formula

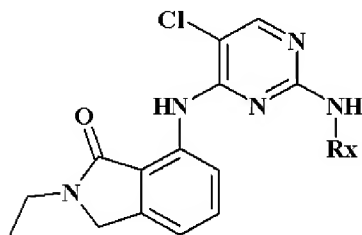


wherein Rx has one of the meanings given in the following table:

Com- pound	Rx
11-1	

11-2	
11-3	
11-4	

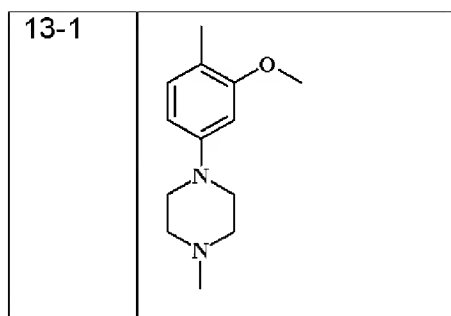
a compound of the formula



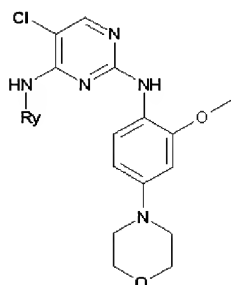
wherein Rx has one of the meanings given in the following table:

Compound.	Rx



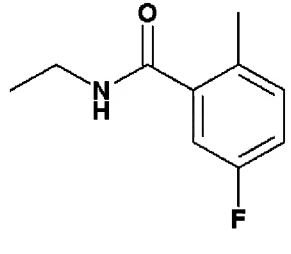
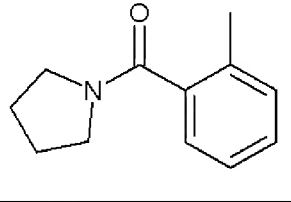
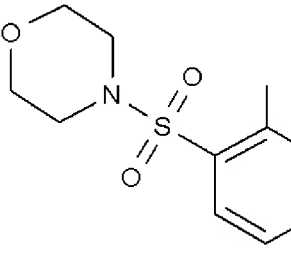


a compound of the formula

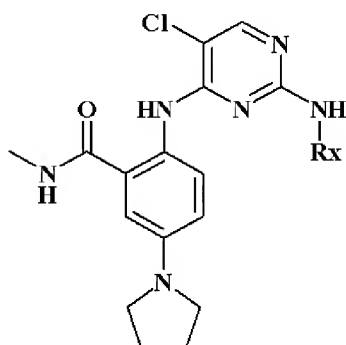


wherein Ry has one of the meanings given in the following table:

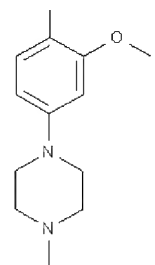
Com- pound	Ry
14-1	
14-2	
14-3	

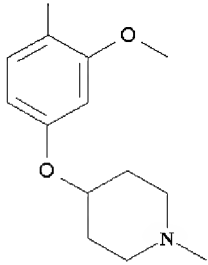
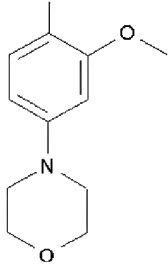
14-5	
14-6	
14-7	

a compound of the formula

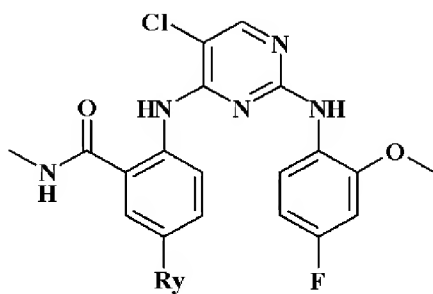


wherein Rx has one of the meanings given in the following table:

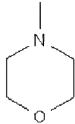
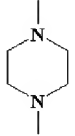

Compound	Rx
15-1	

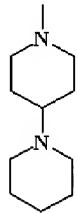
15-2	
15-3	

a compound of the formula

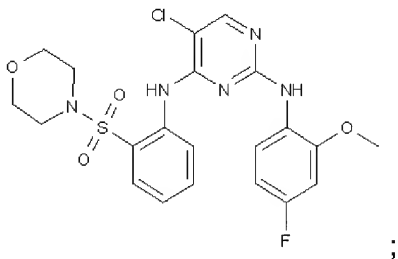


wherein Ry has one of the meanings given in the following table:

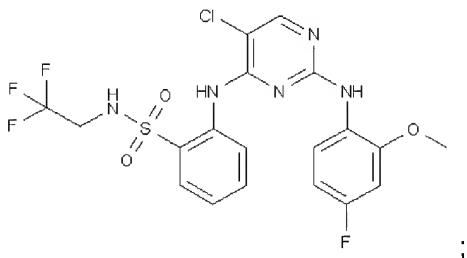
Com- pound	Ry
16-1	
16-2	
16-3	

16-4	
------	---

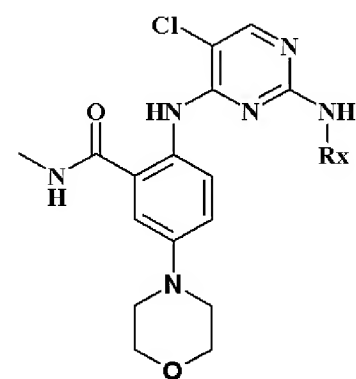
a compound of the formula



a compound of the formula

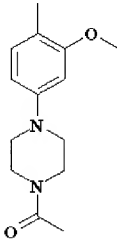


a compound of the formula

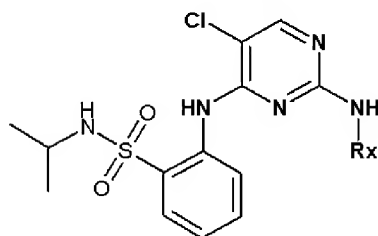


wherein Rx has one of the meanings given in the following table:

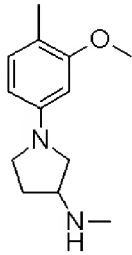
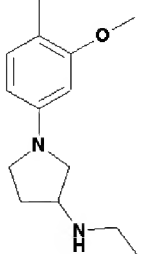
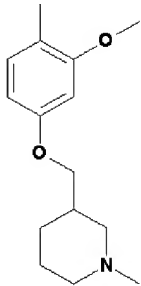
Com- pound.	Rx

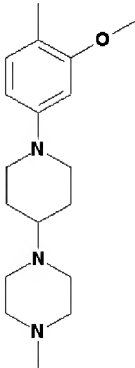
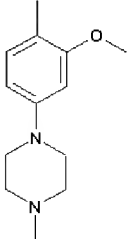
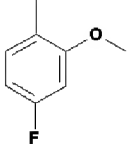
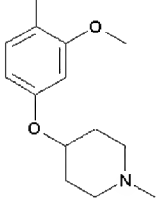
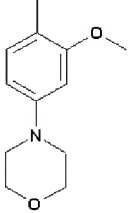
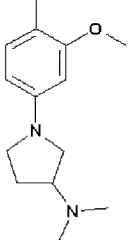
18-1	
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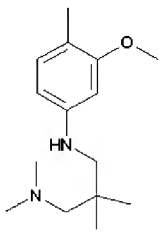
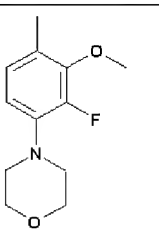
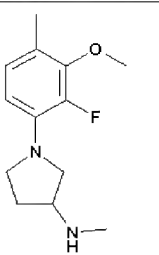
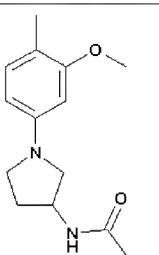
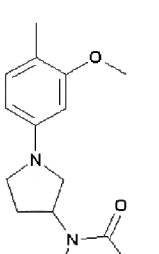
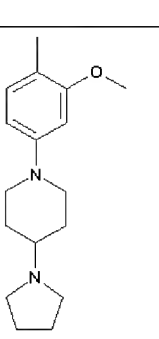
a compound of the formula

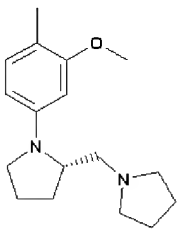
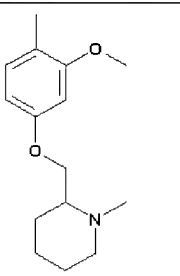
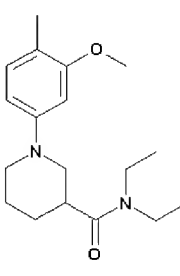
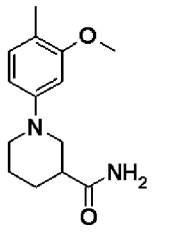
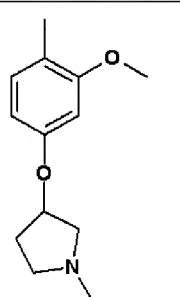


wherein Rx has one of the meanings given in the following table:

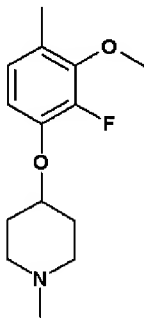
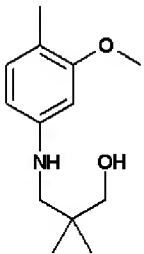
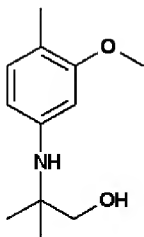
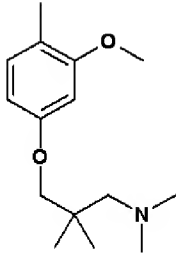
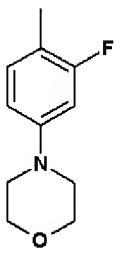
Com- pound	Rx
19-1	
19-2	
19-3	

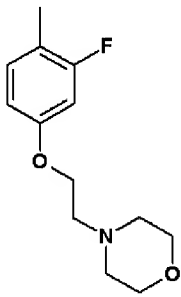
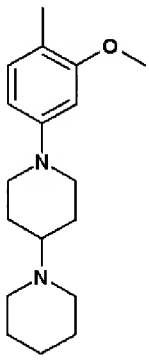
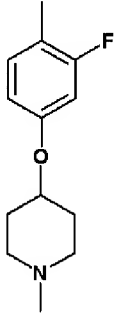
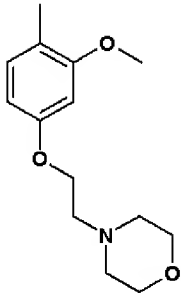
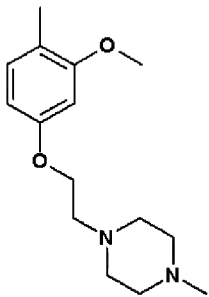
19-4	 <chem>COC1=CC=C(C=C1)N2CCN(CC2)C3CCN(CC3)C4</chem>
19-5	 <chem>COC1=CC=C(C=C1)N2CCN(CC2)C3</chem>
19-6	 <chem>COC1=CC=C(C=C1)N2CCN(CC2)C3</chem>
19-7	 <chem>COC1=CC=C(C=C1)OCC2CCN(C)CC2</chem>
19-8	 <chem>COC1=CC=C(C=C1)N2CCOCC2</chem>
19-9	 <chem>COC1=CC=C(C=C1)N2CCN(C)CC2</chem>

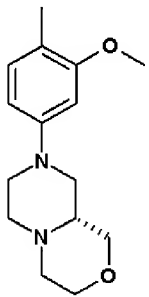
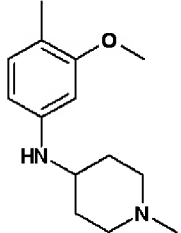
19-10	
19-11	
19-12	
19-13	
19-14	
19-15	

19-16	 <chem>COC1=CC=C(C=C1)N2CC[C@H](C2)N3CCCC3</chem>
19-17	 <chem>COC1=CC=C(C=C1)N2CC[C@H](C2)N3CCCC3</chem>
19-18	 <chem>COC1=CC=C(C=C1)N2CC[C@H](C2)N3CCCC3</chem>
19-19	 <chem>COC1=CC=C(C=C1)N2CC[C@H](C2)N3CCCC3</chem>
19-20	 <chem>COC1=CC=C(C=C1)N2CC[C@H](C2)N3CCCC3</chem>



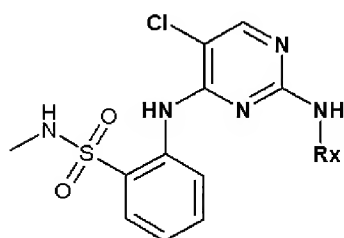
19-21	 <chem>CN1CCN(CC1)Oc2cc(OC)c(F)cc2</chem>
19-22	 <chem>CC(C)(C)C(O)CNc1ccc(OC)cc1</chem>
19-23	 <chem>CC(C)(C)C(O)Nc1ccc(OC)cc1</chem>
19-24	 <chem>CN(C)CC(C)(C)COc1ccc(OC)cc1</chem>
19-25	 <chem>C1CCNCC1c1ccc(F)cc1</chem>

19-26	 <chem>COc1ccc(OCCN2CCCCO2)c(F)c1</chem>
19-27	 <chem>COc1ccc(N2CCCCC2)cc1</chem>
19-28	 <chem>COc1ccc(N2CCCCC2)cc1</chem>
19-29	 <chem>COc1ccc(OCCN2CCCCO2)cc1</chem>
19-30	 <chem>COc1ccc(OCCN(C)CCN(C)C)cc1</chem>

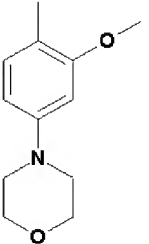
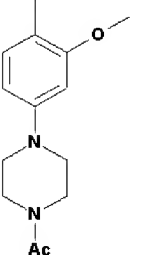
19-31	
19-32	

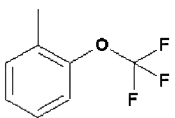
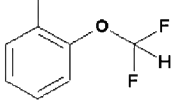
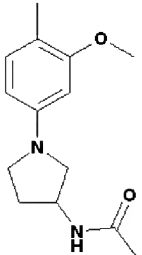
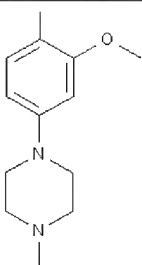
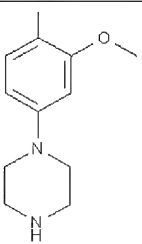
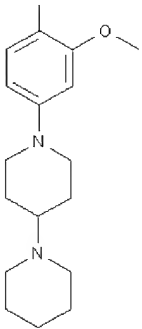
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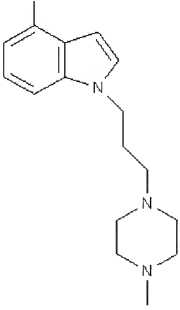
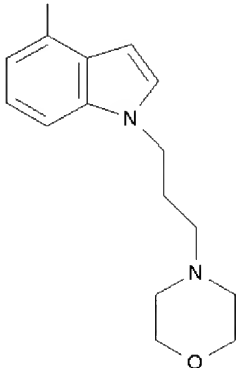
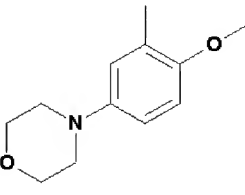
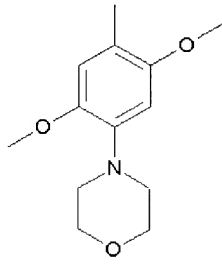
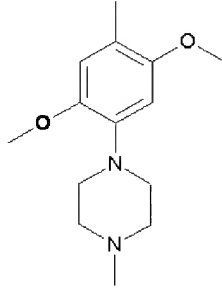
a compound of the formula

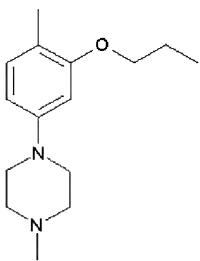
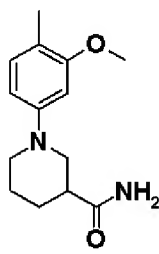
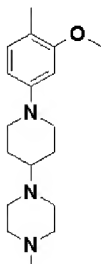
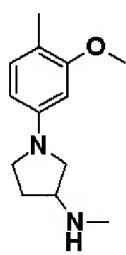
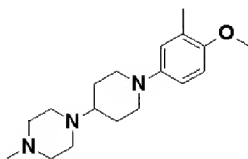
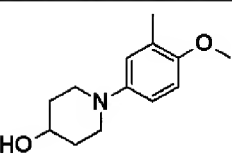
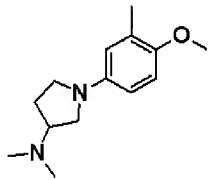


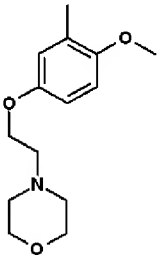
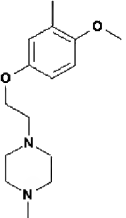
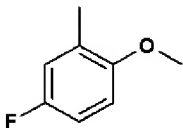
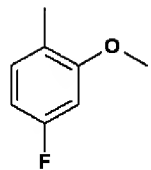
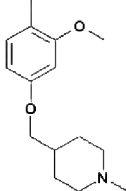
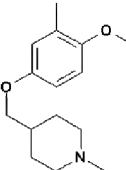
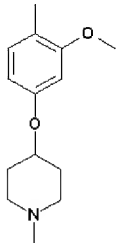
wherein Rx has one of the meanings given in the following table:

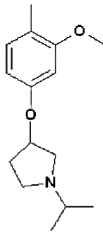
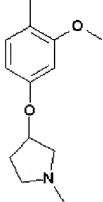
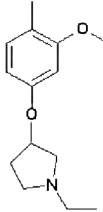
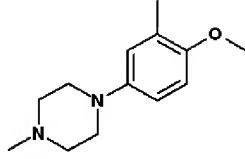
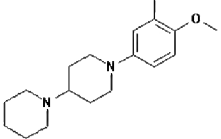
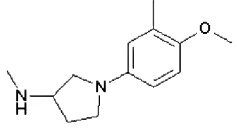
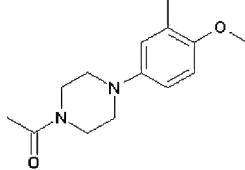
Com- pound	Rx
20-1	
20-2	

20-3	
20-4	
20-5	
20-6	
20-7	
20-8	

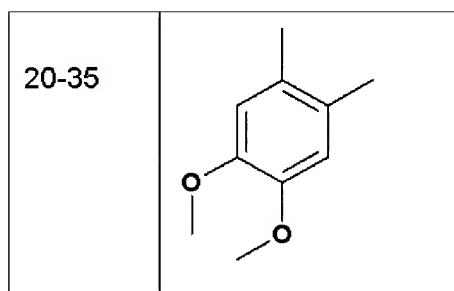
20-9	 <chem>CN1CCN(C)CC1CN2C=CC3=CC=CC=C3C2C</chem>
20-10	 <chem>CN1CCN(C)CC1CN2C=CC3=CC=CC=C3C2C</chem>
20-11	 <chem>CN1CCN(C)CC1CN2C=CC3=CC=CC=C3C2C</chem>
20-12	 <chem>CN1CCN(C)CC1CN2C=CC3=CC=CC=C3C2C</chem>
20-13	 <chem>CN1CCN(C)CC1CN2C=CC3=CC=CC=C3C2C</chem>

20-14	 <chem>CCOC1=CC=C(C=C1)N2CCN(C)CC2</chem>
20-15	 <chem>COc1ccc(cc1)N2CCCCC2C(=O)N</chem>
20-16	 <chem>COc1ccc(cc1)N2CCN(C)CC2</chem>
20-17	 <chem>COc1ccc(cc1)N2CCCC2</chem>
20-18	 <chem>COc1ccc(cc1)N2CCN(C)CC2</chem>
20-19	 <chem>COc1ccc(cc1)N2CCCC(O)C2</chem>
20-20	 <chem>COc1ccc(cc1)N2CCCC2</chem>

20-21	
20-22	
20-23	
20-24	
20-25	
20-26	
20-27	

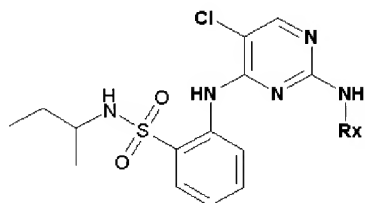
20-28	
20-29	
20-30	
20-31	
20-32	
20-33	
20-34	





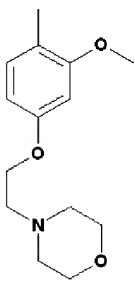
i

a compound of the formula

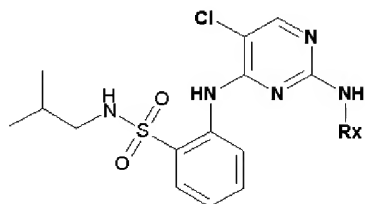


wherein Rx has one of the meanings given in the following table:

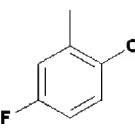
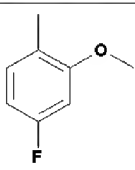
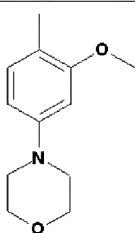
Com- pound	Rx
21-1	
21-2	
21-3	

21-4	
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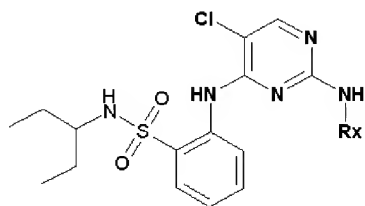
a compound of the formula



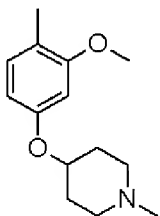
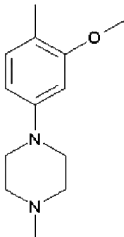
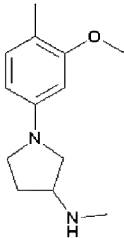
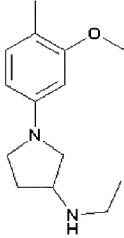
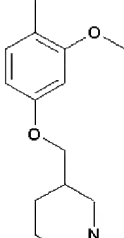
wherein Rx has one of the meanings given in the following table:

Com- pound	Rx
22-1	
22-2	
22-3	

a compound of the formula

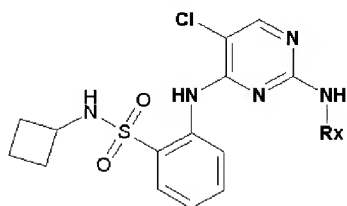


wherein Rx has one of the meanings given in the following table:

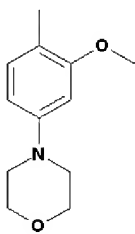
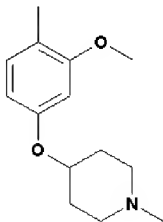
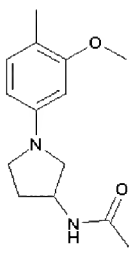
Com- pound	Rx
23-1	
23-2	
23-3	
23-4	
23-5	

23-6	
23-7	
23-8	
23-9	

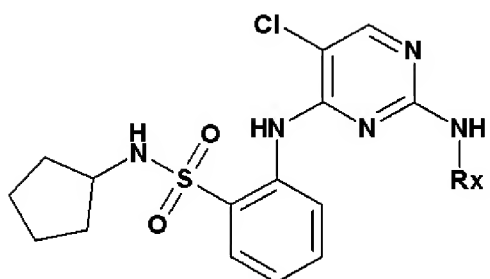
a compound of the formula



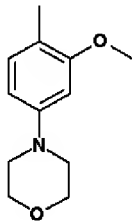
wherein Rx has one of the meanings given in the following table:

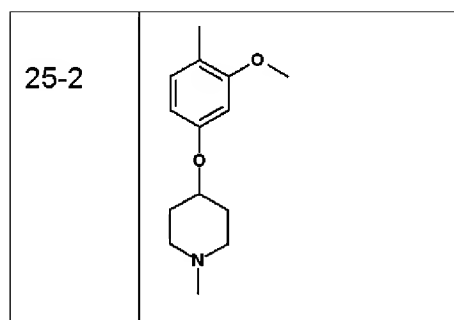
Com- pound	Rx
24-1	
24-2	
24-3	

a compound of the formula

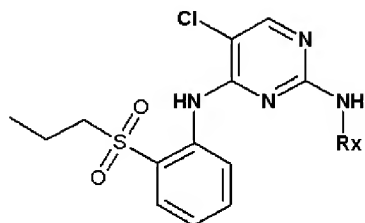


wherein Rx has one of the meanings given in the following table:

Com- pound	Rx
25-1	

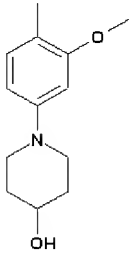
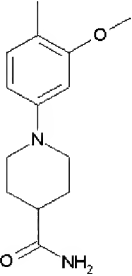
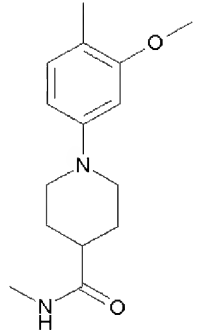
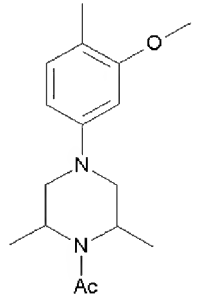
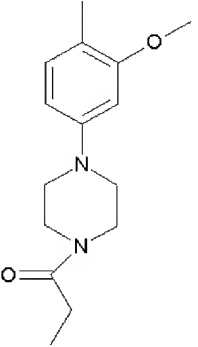


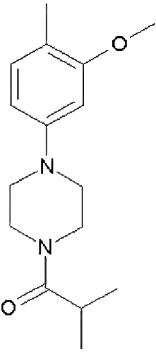
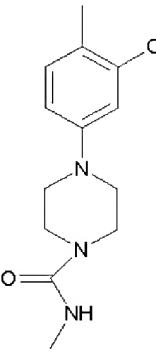
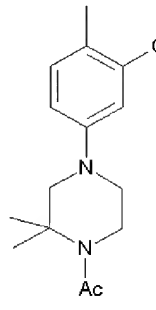
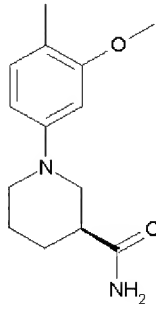
a compound of the formula



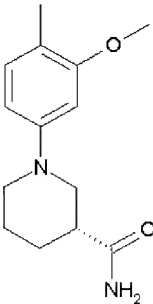
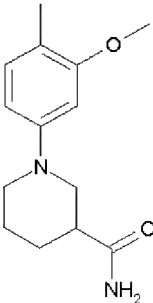
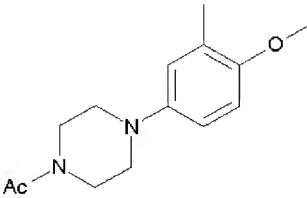
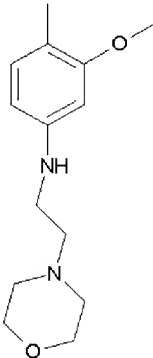
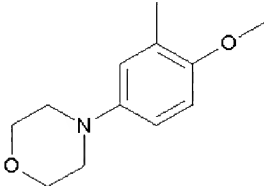
wherein Rx has one of the meanings given in the following table:

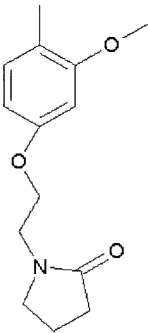
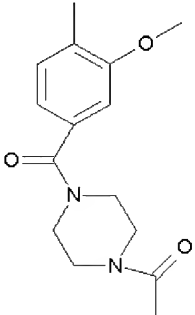
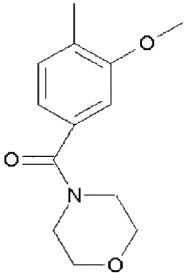
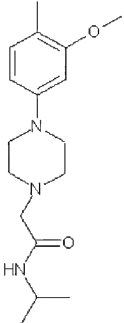
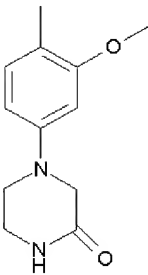
Com- pound	Rx
26-1	
26-2	
26-3	

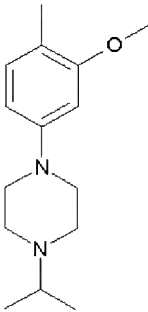
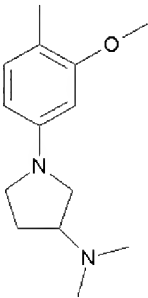
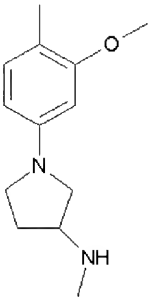
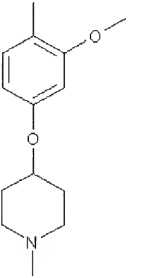
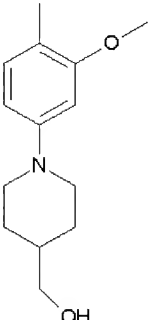
26-4	 <chem>COc1ccc(cc1)N2CCCCC2O</chem>
26-5	 <chem>COc1ccc(cc1)N2CCCCC2C(=O)N</chem>
26-6	 <chem>COc1ccc(cc1)N2CCCCC2C(=O)N(C)C</chem>
26-7	 <chem>COc1ccc(cc1)N2CCN(C)CC2C</chem>
26-8	 <chem>COc1ccc(cc1)N2CCN(CC2)C(=O)CC</chem>

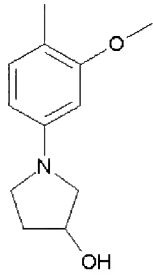
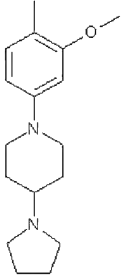
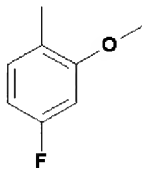
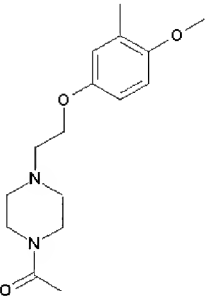
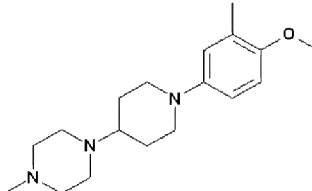
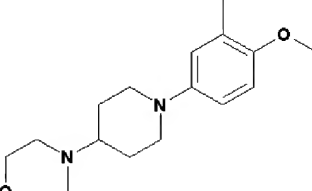
26-9	
26-10	
26-11	
26-12	

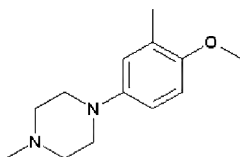
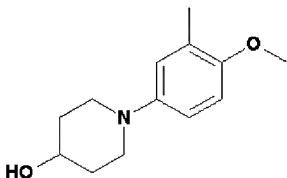
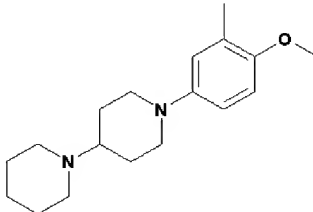
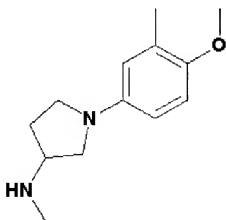
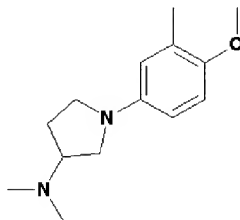
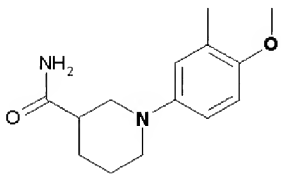


26-13	
26-14	
26-15	
26-16	
26-17	

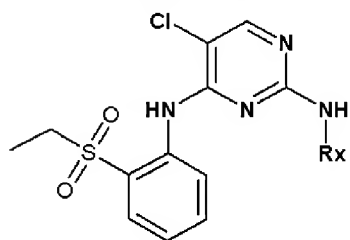
26-18	 <chem>COc1cccc(c1)OCCN2CCCC2=O</chem>
26-19	 <chem>COc1cccc(c1)C(=O)N1CCN(C)CC1</chem>
26-20	 <chem>COc1cccc(c1)C(=O)N1CCOCC1</chem>
26-21	 <chem>COc1cccc(c1)N1CCN(CC1)CC(=O)N(C)C</chem>
26-22	 <chem>COc1cccc(c1)N1CCNC(=O)CC1=O</chem>

26-23	 <chem>CN1CCN(CC1)c2ccc(OC)cc2</chem>
26-24	 <chem>CN1CCN(C1)c2ccc(OC)cc2</chem>
26-25	 <chem>C1CCN(C1)c2ccc(OC)cc2</chem>
26-26	 <chem>C1CCN(CC1)c2ccc(OC)cc2</chem>
26-27	 <chem>C1CCN(CC1CO)c2ccc(OC)cc2</chem>

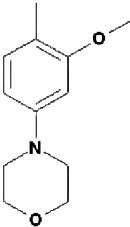
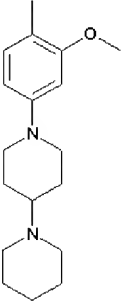
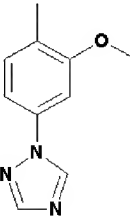
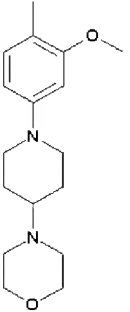
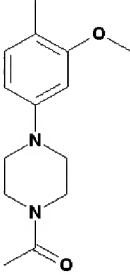
26-28	
26-29	
26-30	
26-31	
26-32	
26-33	

26-34	
26-35	
26-36	
26-37	
26-38	
26-39	

a compound of the formula

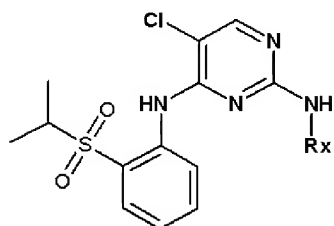


wherein Rx has one of the meanings given in the following table:

Compound	Rx
27-1	
27-2	
27-3	
27-4	
27-5	

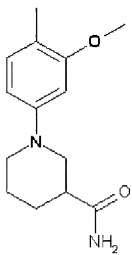
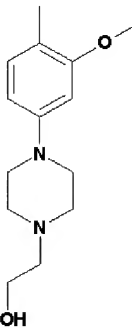
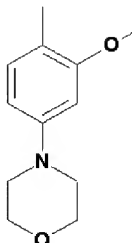
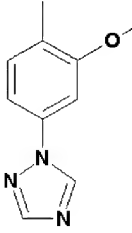
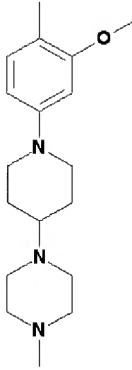
27-6	
27-7	
27-8	
27-9	

a compound of the formula

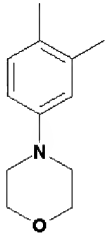
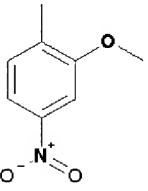
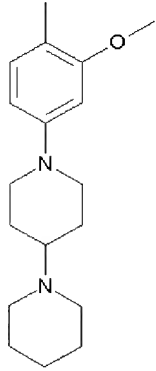
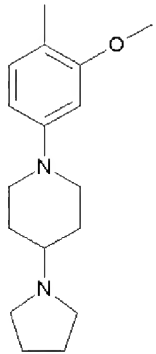
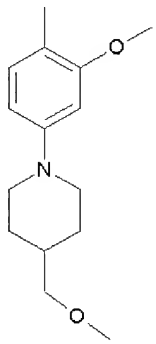


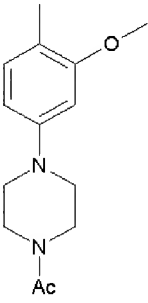
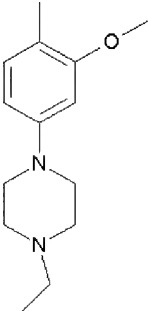
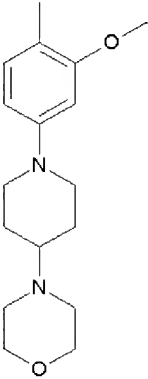
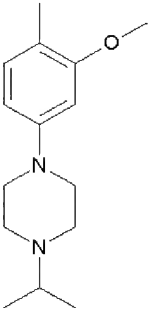
wherein Rx has one of the meanings given in the following table:

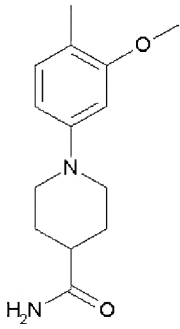
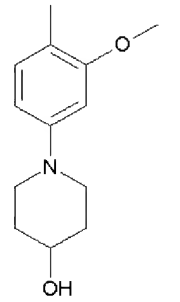
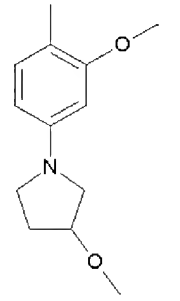
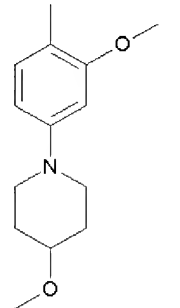
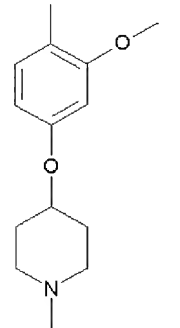
Com- pound	Rx

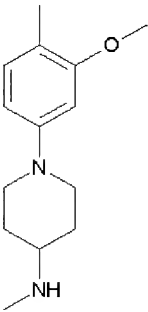
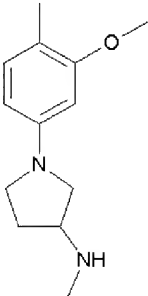
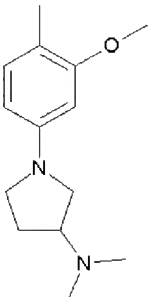
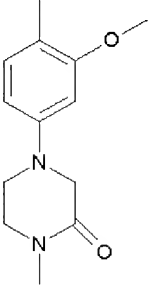
28-1	
28-2	
28-3	
28-4	
28-5	

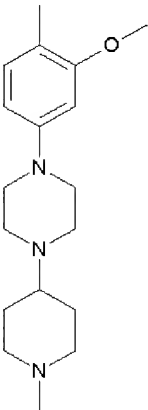
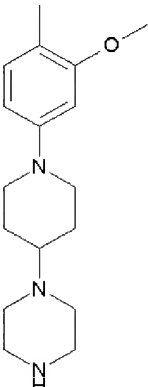
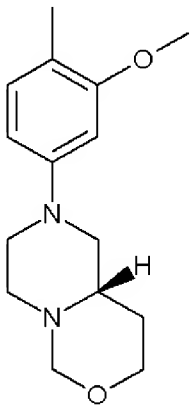
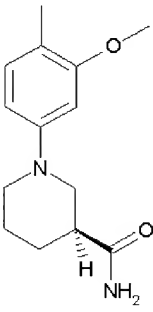


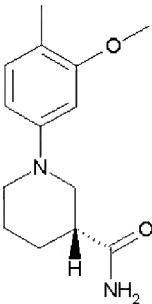
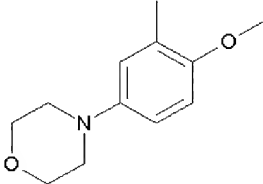
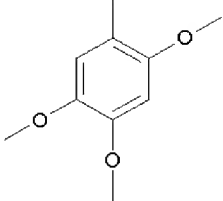
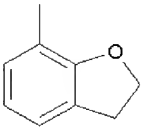
28-6	 <chem>Cc1cc(C)cc(N2CCCC2)c1</chem>
28-7	 <chem>COc1cc([N+](=O)[O-])ccc1N2CCCC2</chem>
28-8	 <chem>COc1cc(ccc1N2CCCCC2)N3CCCCC3</chem>
28-9	 <chem>COc1cc(ccc1N2CCCCC2)N3CCCC3</chem>
28-10	 <chem>COc1cc(ccc1N2CCCCC2)COC3CCCCC3</chem>

28-11	
28-12	
28-13	
28-14	

28-15	
28-16	
28-17	
28-18	
28-19	

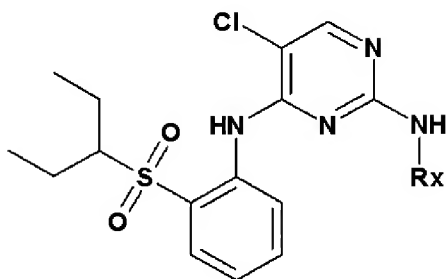
28-20	 <chem>CN1CCN(CC1)c2ccc(OC)cc2</chem>
28-21	 <chem>CN1CCN(C1)c2ccc(OC)cc2</chem>
28-22	 <chem>CN(C)C1CCN(C1)c2ccc(OC)cc2</chem>
28-23	 <chem>CN1CC(=O)N(C)CC1Nc2ccc(OC)cc2</chem>

28-24	 <chem>COc1ccc(cc1)N2CCNCC2</chem>
28-25	 <chem>COc1ccc(cc1)N2CCCCC2</chem>
28-26	 <chem>COc1ccc(cc1)N2CCN(C2)C3CCOCC3</chem>
28-27	 <chem>COc1ccc(cc1)N2CCCC[C@H]2C(=O)N</chem>

28-28	
28-29	
28-30	
28-31	

;

a compound of the formula



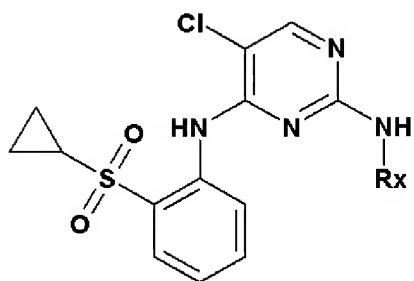
wherein Rx has one of the meanings given in the following table:

Com- pound	Rx

29-1	
29-2	
29-3	

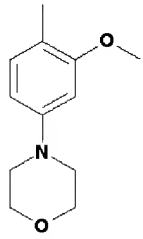
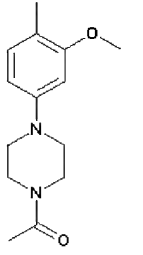
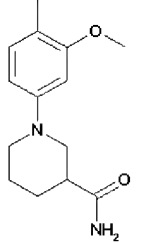
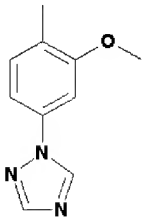
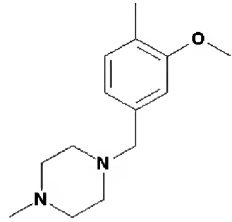
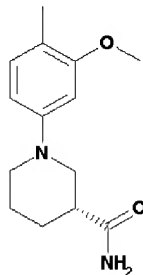
;

a compound of the formula

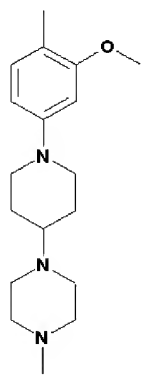


wherein Rx has one of the meanings given in the following table:

Com- pound	Rx
---------------	----

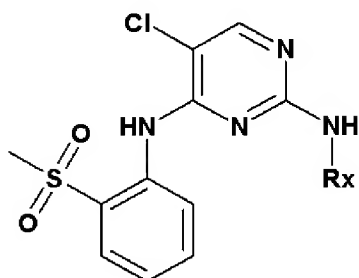
30-1	
30-2	
30-3	
30-4	
30-5	
30-6	



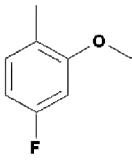
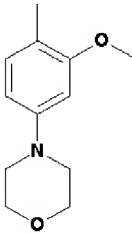
30-7	
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;

a compound of the formula

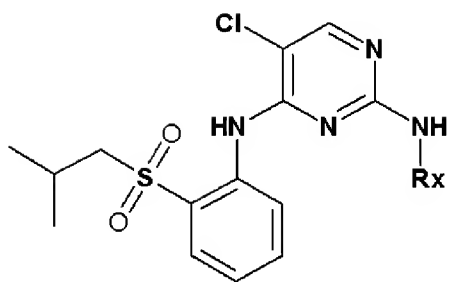


wherein Rx has one of the meanings given in the following table:

Com- pound	Rx
31-1	
31-2	

;

a compound of the formula

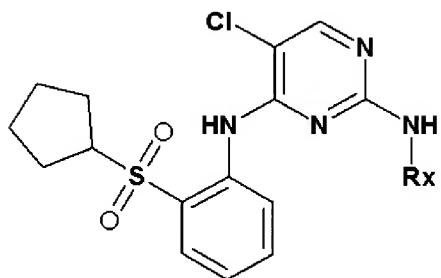


wherein Rx has one of the meanings given in the following table:

Com- pounds	Rx
32-1	
32-2	

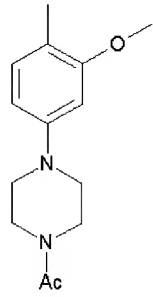
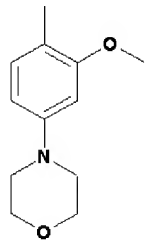
:

a compound of the formula



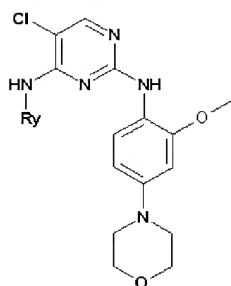
wherein Rx has one of the meanings given in the following table:

Com- pound	Rx
---------------	----

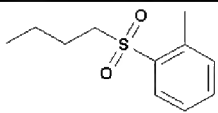
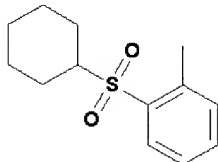
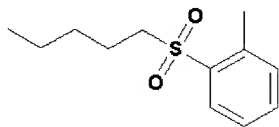
33-1	
33-2	

;

a compound of the formula



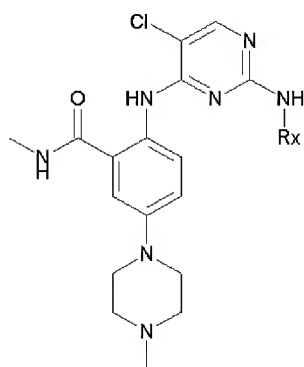
wherein Ry has one of the meanings given in the following table:

Com- pound	Ry
34-1	
34-2	
34-3	

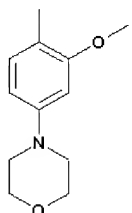
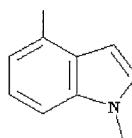
34-4	
34-5	
34-6	
34-7	
34-8	

;

a compound of the formula



wherein Rx has one of the meanings given in the following table:

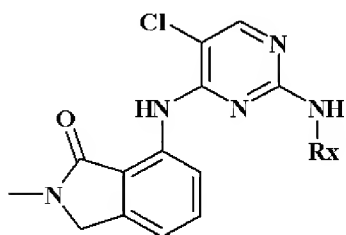
35-1	
35-2	

2-[5-bromo-2-(2-methoxy-4-morpholin-4-yl-phenylamino)-pyrimidin-4-ylamino]-N, N-dimethyl-benzenesulfonamide;

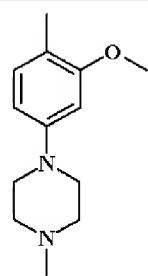
2-[5-bromo-2-(2-methoxy-4-morpholin-4-yl-phenylamino)-pyrimidin-4-ylamino]-5-fluoro-N-methyl-benzenesulfonamide;

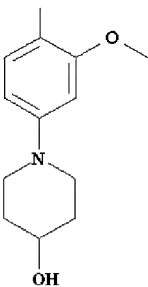
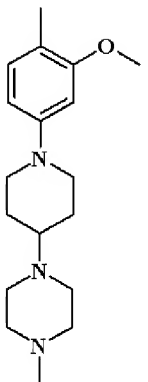
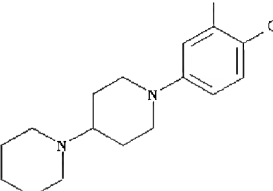
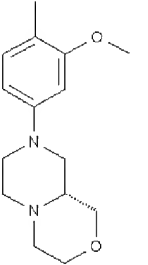
7-[5-chloro-2-(2-methoxy-4-morpholin-4-yl-phenylamino)-pyrimidin-4-ylamino]-2-methyl-2,3-dihydro-isoindol-1-one;

and a compound of the formula



wherein Rx has one of the meanings given in the following table:

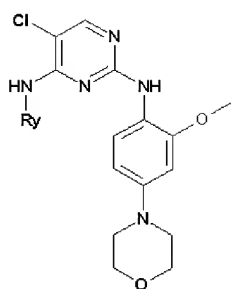
Compound	Rx
12-2	

12-3	
12-4	
12-5	
12-6	

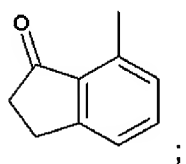
;

or a pharmaceutically acceptable salt thereof.

**24. (New)** A compound of the formula



wherein Ry has the formula



or a pharmaceutically acceptable salt thereof.